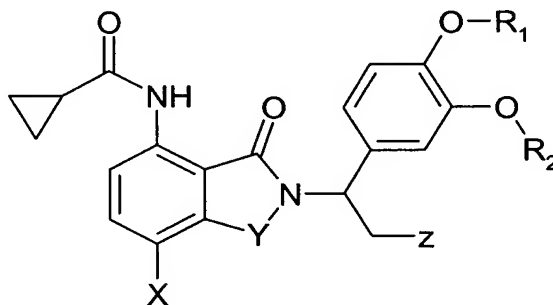


IN THE CLAIMS:

The listing of claims below will replace all prior versions and listings of claims in this application.

Listing Of Claims:

1. (Currently amended) A compound of formula (I):



wherein:

Y is -C(O)-, -CH₂-, -CH₂C(O)- or -SO₂-;

X is H;

Z is (C₀₋₄-alkyl)-C(O)R³, C₁₋₄-alkyl, (C₀₋₄-alkyl)-OH, (C₁₋₄-alkyl)-O-(C₁₋₄-alkyl), (C₁₋₄-alkyl)-SO₂(C₁₋₄-alkyl), (C₀₋₄-alkyl)-SO(C₁₋₄-alkyl), (C₀₋₄-alkyl)-NH₂, (C₀₋₄-alkyl)-N(C₁₋₈-alkyl)₂, (C₀₋₄-alkyl)-N(H)(OH), ~~(C₀₋₄-alkyl)-dichloropyridine~~, or CH₂NSO₂-(C₁₋₄-alkyl);

R₁ and R₂ are independently C₁₋₈-alkyl, cycloalkyl, or (C₁₋₄-alkyl)-cycloalkyl;

R³ is, NR⁴ R⁵, OH, or O-(C₁₋₈-alkyl);

R⁴ is H;

R⁵ is -OH, or -O-C(O)R⁶;

R⁶ is C₁₋₈-alkyl, amino-(C₁₋₈-alkyl), (C₁₋₈-alkyl)-(C₃₋₆-cycloalkyl), C₃₋₆-cycloalkyl, phenyl, benzyl, or aryl;

or a pharmaceutically acceptable salt, ~~or~~ solvate, stereoisomer or clathrate thereof.

2. (Original) The compound of claim 1, wherein Z is (C₀₋₄-alkyl)-C(O)R³, C₁₋₄-alkyl, (C₀₋₄-alkyl)-OH, (C₁₋₄-alkyl)-O(C₁₋₄-alkyl), (C₁₋₄-alkyl)-SO₂(C₁₋₄-alkyl), (C₀₋₄-alkyl)-SO(C₁₋₄-alkyl), (C₀₋₄-alkyl)-NH₂, (C₀₋₄-alkyl)-N(C₁₋₈-alkyl)₂, (C₀₋₄-alkyl)-N(H)(OH), or CH₂NSO₂(C₁₋₄-alkyl).

3-4. (Canceled)

5. (Currently amended) The compound of claim 1 ~~or 3~~ wherein R₁ is CH₃.

6. (Currently amended) The compound of claim 1 ~~or 3~~ wherein R₂ is CH₂CH₃, CH₃, CH₂-cyclopropyl, or cyclopentyl.

7. (Currently amended) The compound of claim 1 ~~or 3~~ wherein Y is -C(O)- or -CH₂-.

8-10. (Canceled)

11. (Currently amended) An enantiomerically pure S isomer of a compound of claim 1, ~~3 or 9~~, substantially free of its R isomer, or a pharmaceutically acceptable salt, solvate, ~~hydrate~~, stereoisomer, or clathrate, ~~or prodrug~~ thereof.

12. (Currently amended) An enantiomerically pure R isomer of a compound of claim 1, ~~3 or 9~~, substantially free of its S isomer, or a pharmaceutically acceptable salt, solvate, ~~hydrate~~, stereoisomer, or clathrate, ~~or prodrug~~ thereof.

13. (Currently amended) A compound, wherein the compound is:
(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-3-hydroxy-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;
(3R)-(tert-Butoxy)-N-{3-[7-(cyclopropylcarbonylamino)-1-oxoisindolin-2-yl]-3-(3-ethoxy-4-methoxyphenyl)propyl}carbonylamino (tert-butoxy)formate;
(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-3-hydroxyamino-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;
(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-3-methanesulfonylamino-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;
(1R)-Cyclopropanecarboxylic acid {2-[3-amino-1-(3-ethoxy-4-methoxy-phenyl)-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;
(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-3-ureido-propyl]-3-oxo-

2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[3-dimethylamino-1-(3-ethoxy-4-methoxy-phenyl)-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide hydrochloride;

(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-3-methanesulfonyl-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-2-hydroxycarbamoyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[2-acetoxycarbamoyl-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(3R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-3-methanesulfinyl-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(3R)-3-[4-Chloro-7-(cyclopropanecarbonyl-amino)-1-oxo-1,3-dihydro-isoindol-2-yl]-3-(3-ethoxy-4-methoxy-phenyl)-propionic acid;

(3R)-3-[4-Chloro-7-(cyclopropanecarbonyl-amino)-1-oxo-1,3-dihydro-isoindol-2-yl]-3-(3-ethoxy-4-methoxy-phenyl)-propionic acid methyl ester;

(1R)-Cyclopropanecarboxylic acid {2-[2-carbamoyl-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-7-chloro-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {7-chloro-2-[2-dimethylcarbamoyl-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {7-chloro-2-[1-(3-ethoxy-4-methoxy-phenyl)-2-hydroxycarbamoyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[2-acetoxycarbamoyl-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-7-chloro-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1S)-Cyclopropanecarboxylic acid {7-chloro-2-[1-(3-ethoxy-4-methoxy-phenyl)-2-methanesulfonyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1S)-Cyclopropanecarboxylic acid {7-bromo-2-[1-(3-ethoxy-4-methoxy-phenyl)-2-methanesulfonyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-propyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

~~Cyclopropanecarboxylic acid {2-[2-(3,5-dichloro-pyridin-4-yl)-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;~~

(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-3-hydroxy-3-methyl-

butyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[2-cyclopropanecarbonyloxycarbamoyl-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-2-isobutyryloxycarbamoyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[2-(2,2-dimethyl-propionyloxycarbamoyl)-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1R)-Cyclopropanecarboxylic acid {2-[2-(3,3-dimethyl-butyryloxycarbamoyl)-1-(3-ethoxy-4-methoxy-phenyl)-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

(1S)-Cyclopropanecarboxylic acid {2-[1-(3-ethoxy-4-methoxy-phenyl)-2-methanesulfonyl-ethyl]-7-fluoro-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-amide;

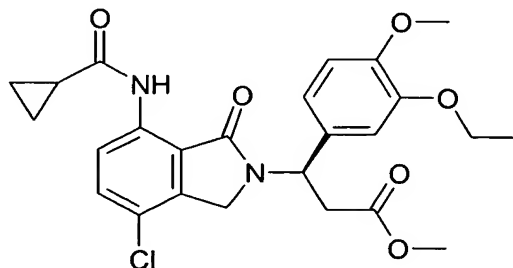
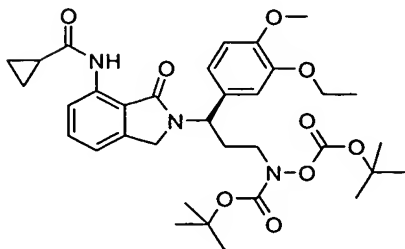
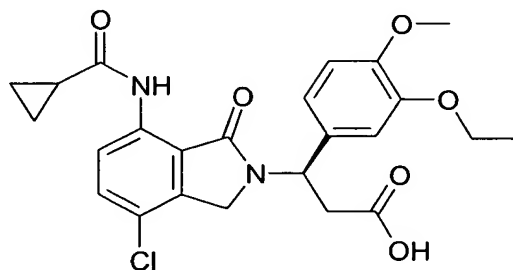
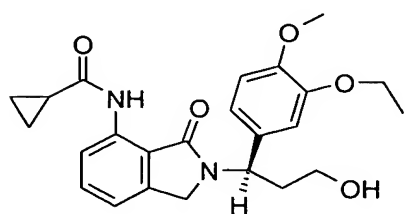
(1S)-3-{7-Chloro-2-[1-(3-ethoxy-4-methoxy-phenyl)-2-methanesulfonyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-1,1-dimethyl-urea;

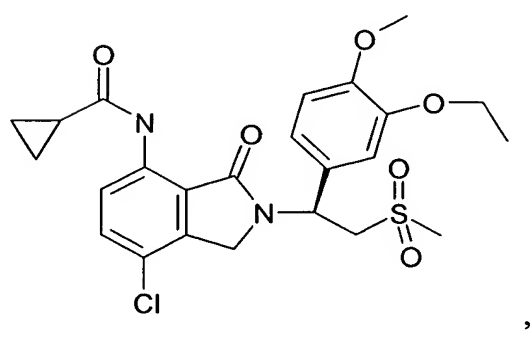
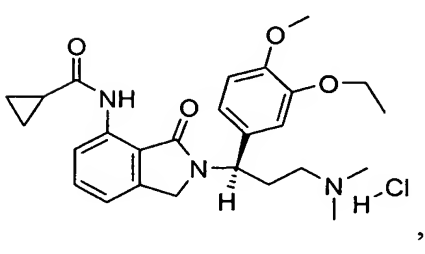
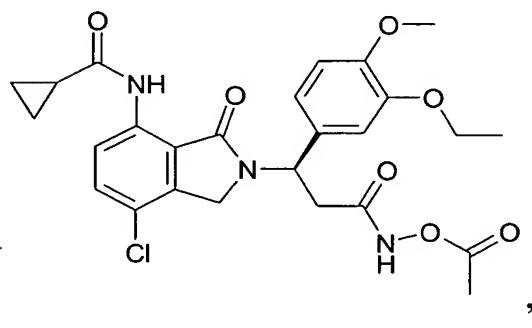
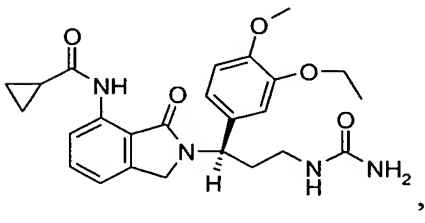
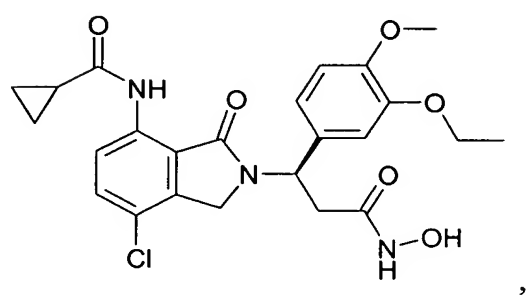
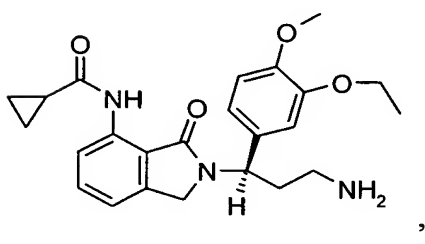
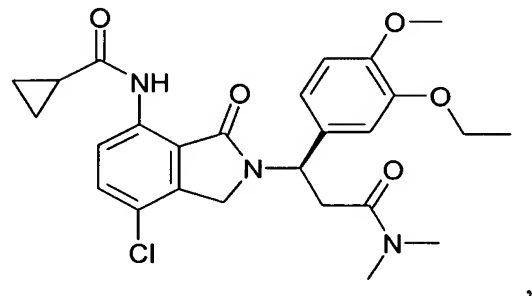
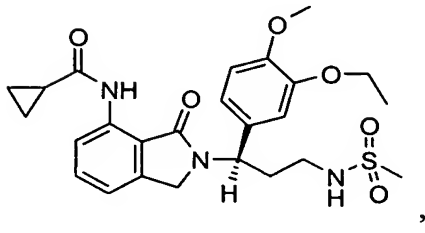
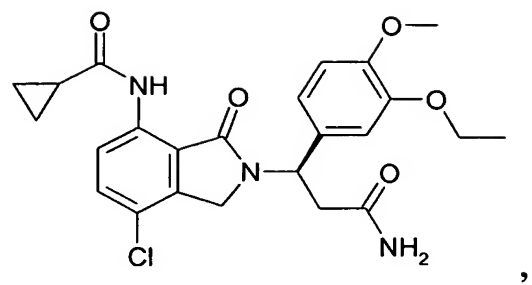
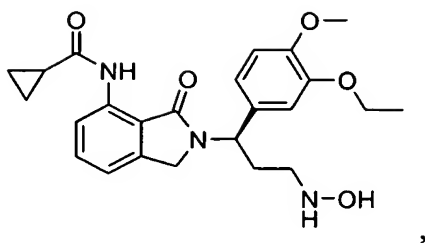
(1S)-N-{7-Chloro-2-[1-(3-ethoxy-4-methoxy-phenyl)-2-methanesulfonyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-2-(4-methyl-piperazin-1-yl)-acetamide;

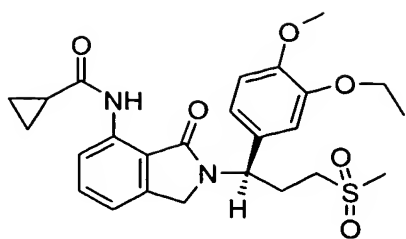
(1S)-N-{7-Chloro-2-[1-(3-ethoxy-4-methoxy-phenyl)-2-methanesulfonyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-2-morpholin-4-yl-acetamide hydrochloride;

(1S)-N-{7-Chloro-2-[1-(3-ethoxy-4-methoxy-phenyl)-2-methanesulfonyl-ethyl]-3-oxo-2,3-dihydro-1H-isoindol-4-yl}-2-dimethylamino-acetamide hydrochloride.

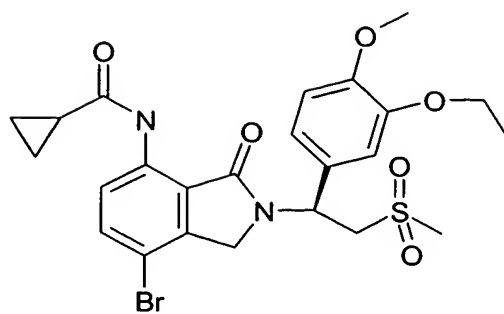
14. (Currently amended) A compound of the formula:



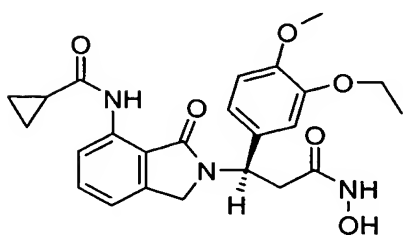




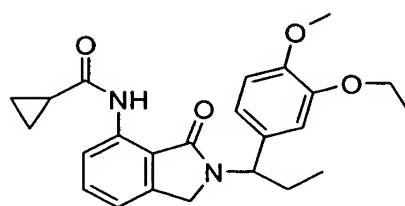
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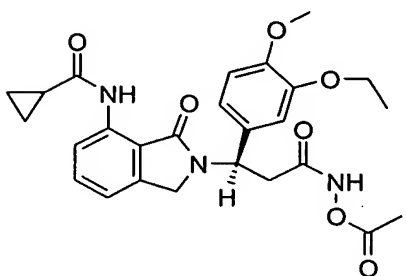
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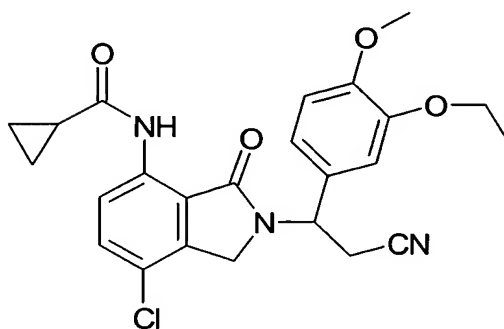
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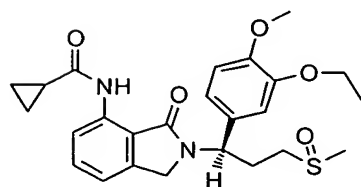
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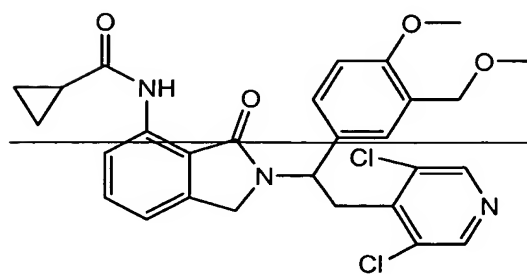
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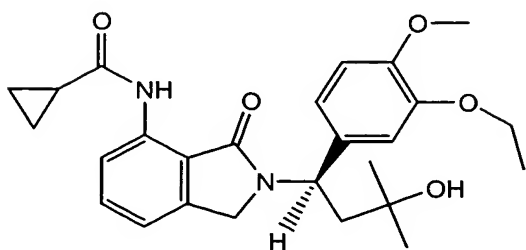
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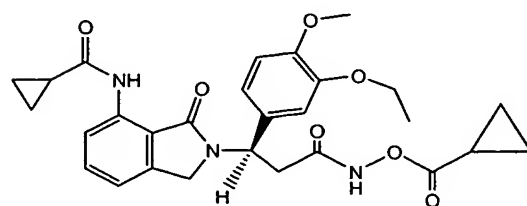
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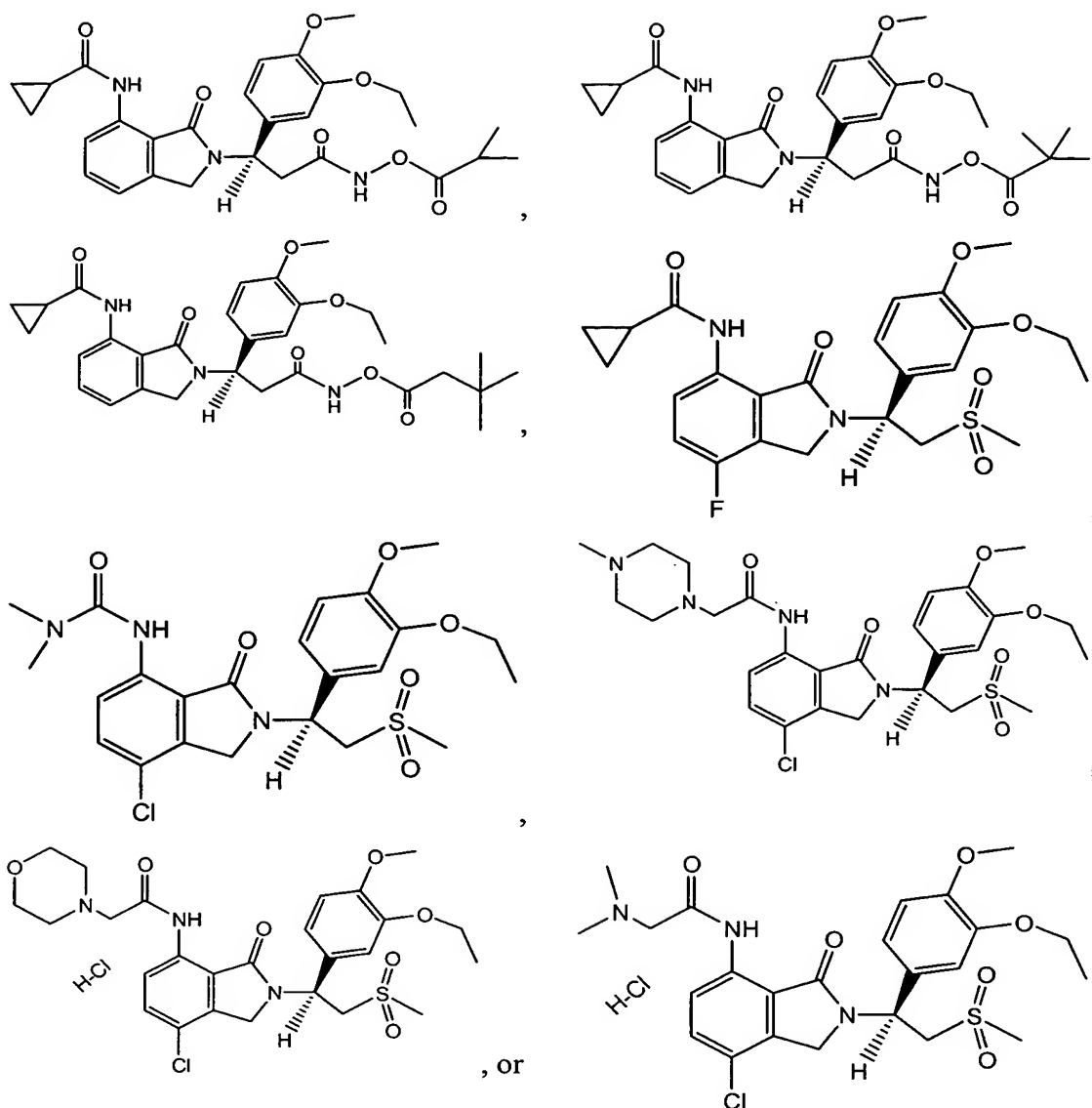
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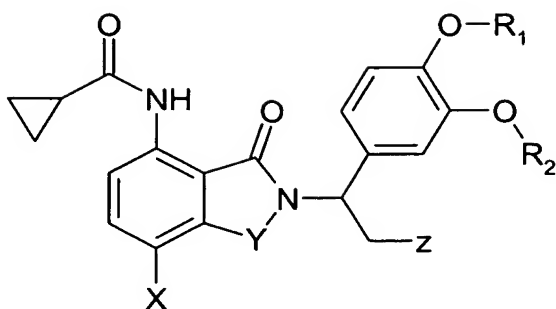
,



or a pharmaceutically acceptable salt or solvate thereof.

15. (Canceled)

16. (Currently amended) A pharmaceutical composition comprising a pharmaceutically acceptable carrier, excipient, or diluent and a compound of formula (I):



wherein:

Y is -C(O)-, -CH₂-, -CH₂C(O)- or -SO₂-;

X is H;

Z is (C₀₋₄-alkyl)-C(O)R³, C₁₋₄-alkyl, (C₀₋₄-alkyl)-OH, (C₁₋₄-alkyl)-O-(C₁₋₄-alkyl), (C₁₋₄-alkyl)-SO₂(C₁₋₄-alkyl), (C₀₋₄-alkyl)-SO(C₁₋₄-alkyl), (C₀₋₄-alkyl)-NH₂, (C₀₋₄-alkyl)-N(C₁₋₈-alkyl)₂, (C₀₋₄-alkyl)-N(H)(OH), ~~(C₀₋₄-alkyl)-dichloropyridine~~, or CH₂NSO₂-(C₁₋₄-alkyl);

R₁ and R₂ are independently C₁₋₈-alkyl, cycloalkyl, or (C₁₋₄-alkyl)-cycloalkyl;

R³ is, NR⁴ R⁵, OH, or O-(C₁₋₈-alkyl);

R⁴ is H;

R⁵ is -OH, or -O-C(O)R⁶;

R⁶ is C₁₋₈-alkyl, amino-(C₁₋₈-alkyl), (C₁₋₈-alkyl)-(C₃₋₆-cycloalkyl), C₃₋₆-cycloalkyl, phenyl, benzyl, or aryl;

or a pharmaceutically acceptable salt or solvate thereof.

17-21. (Canceled)

22. (New) The compound of claim 1, wherein said solvate is a hydrate.

23. (New) The enantiomerically pure S isomer of claim 11, wherein said solvate is a hydrate.

24. (New) The enantiomerically pure R isomer of claim 12, , wherein said solvate is a hydrate.